

```
ring nodes:
    1 2 3 4 5
ring/chain nodes:
    9 10
chain bonds:
    1-6 2-9 6-7 7-10
ring bonds:
    1-2 1-5 2-3 3-4 4-5
exact/norm bonds:
    1-2 1-6 2-3 2-9 6-7 7-10
exact bonds:
    1-5 3-4 4-5
isolated ring systems:
    containing 1:

Match level:
    1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:Atom 9:CLASS 10:CLASS 11:CLASS
```

NEWS 37

NEWS 38

NEWS 39

Dec 17

Dec 30

Jan 13

Welcome to STN International! Enter x:x

LOGINID:ssspta1611txm PASSWORD: TERMINAL (ENTER 1, 2, 3, OR ?):2 Welcome to STN International NEWS Web Page URLs for STN Seminar Schedule - N. America "Ask CAS" for self-help around the clock Apr 08 NEWS 2 NEWS 3 Apr 09 BEILSTEIN: Reload and Implementation of a New Subject Area NEWS 4 Apr 09 ZDB will be removed from STN NEWS 5 Apr 19 US Patent Applications available in IFICDB, IFIPAT, and IFIUDB NEWS 6 Apr 22 Records from IP.com available in CAPLUS, HCAPLUS, and ZCAPLUS NEWS Apr 22 BIOSIS Gene Names now available in TOXCENTER NEWS Apr 22 Federal Research in Progress (FEDRIP) now available NEWS 9 Jun 03 New e-mail delivery for search results now available NEWS 10 Jun 10 MEDLINE Reload NEWS 11 Jun 10 PCTFULL has been reloaded NEWS 12 Jul 02 FOREGE no longer contains STANDARDS file segment NEWS 13 Jul 22 USAN to be reloaded July 28, 2002; saved answer sets no longer valid Jul 29 NEWS 14 Enhanced polymer searching in REGISTRY NEWS 15 Jul 30 NETFIRST to be removed from STN NEWS 16 CANCERLIT reload Aug 08 NEWS 17 Aug 08 PHARMAMarketLetter(PHARMAML) - new on STN NEWS 18 Aug 08 NTIS has been reloaded and enhanced NEWS 19 Aug 19 Aquatic Toxicity Information Retrieval (AQUIRE) now available on STN NEWS 20 Aug 19 IFIPAT, IFICDB, and IFIUDB have been reloaded NEWS 21 The MEDLINE file segment of TOXCENTER has been reloaded Aug 19 Sequence searching in REGISTRY enhanced NEWS 22 Aug 26 NEWS 23 Sep 03 JAPIO has been reloaded and enhanced NEWS 24 Sep 16 Experimental properties added to the REGISTRY file NEWS 25 Sep 16 CA Section Thesaurus available in CAPLUS and CA NEWS 26 Oct 01 CASREACT Enriched with Reactions from 1907 to 1985 NEWS 27 Oct 21 EVENTLINE has been reloaded NEWS 28 Oct 24 BEILSTEIN adds new search fields NEWS 29 Oct 24 Nutraceuticals International (NUTRACEUT) now available on STN NEWS 30 Oct 25 MEDLINE SDI run of October 8, 2002 NEWS 31 Nov 18 DKILIT has been renamed APOLLIT NEWS 32 Nov 25 More calculated properties added to REGISTRY NEWS 33 Dec 02 TIBKAT will be removed from STN NEWS 34 Dec 04 CSA files on STN PCTFULL now covers WP/PCT Applications from 1978 to date NEWS 35 Dec 17 NEWS 36 Dec 17 TOXCENTER enhanced with additional content

NEWS EXPRESS January 6 CURRENT WINDOWS VERSION IS V6.01a,
CURRENT MACINTOSH VERSION IS V6.0b(ENG) AND V6.0Jb(JP),
AND CURRENT DISCOVER FILE IS DATED 01 OCTOBER 2002
NEWS HOURS STN Operating Hours Plus Help Desk Availability
NEWS INTER General Internet Information

ISMEC no longer available

Adis Clinical Trials Insight now available on STN

Indexing added to some pre-1967 records in CA/CAPLUS

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CAS World Wide Web Site (general information) NEWS WWW

Enter NEWS followed by the item number or name to see news on that specific topic.

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\* \* \* \* \* \* \* \* STN Columbus

FILE 'HOME' ENTERED AT 19:03:27 ON 15 JAN 2003

=> file req

COST IN U.S. DOLLARS

SINCE FILE ENTRY SESSION 0.21

TOTAL

FULL ESTIMATED COST

FILE 'REGISTRY' ENTERED AT 19:03:34 ON 15 JAN 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 14 JAN 2003 HIGHEST RN 479024-64-1 DICTIONARY FILE UPDATES: 14 JAN 2003 HIGHEST RN 479024-64-1

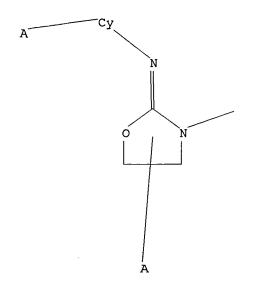
TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

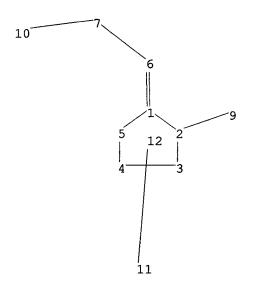
Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

Uploading C:\Program Files\Stnexp\Queries\10004396.str





chain nodes : 6 7 11 ring nodes : 1 2 3 4 5 ring/chain nodes : 9 10 chain bonds : 1-6 2-9 6-7 7-10 ring bonds : 1-2 1-5 2-3 3-4 4-5 exact/norm bonds : 1-2 1-6 2-3 2-9 6-7 7-10 exact bonds : 1-5 3-4 4-5 isolated ring systems : containing 1 :

Match level:
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:Atom 9:CLASS 10:CLASS 11:CLASS 12:CLASS

### L1 STRUCTURE UPLOADED

=> s l1 SAMPLE SEARCH INITIATED 19:03:50 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 101 TO ITERATE

100.0% PROCESSED 101 ITERATIONS 15 ANSWERS SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*
PROJECTED ITERATIONS: 1418 TO 2622
PROJECTED ANSWERS: 68 TO 532

L2 15 SEA SSS SAM L1

=> d scan

L2 15 ANSWERS REGISTRY COPYRIGHT 2003 ACS IN Benzenamine, N-[4-(chloromethyl)-3-phenyl-2-oxazolidinylidene]- (9CI) MF C16 H15 C1 N2 O

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT \*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1): Uploading 'Upload SSTN' IS NOT VALID HERE

To display more answers, enter the number of answers you would like to see. To end the display, enter "NONE", "N", "0", or "END". HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):C:\Program Files\Stnexp\Queries\10004396.str

To display more answers, enter the number of answers you would like to see. To end the display, enter "NoNE", "N", "0", or "END". How MANY MORE ANSWERS DO YOU WISH TO SCAN? (1): "0 SZ'  $\theta$ - $\theta$ - $\theta$ -1" J\*' IS NOT VALID HERE

To display more answers, enter the number of answers you would like to see. To end the display, enter "NONE", "N", "0", or "END". HOW MANY MORE ANSWERS DO YOU WISH TO SCAM? (1): "0 S2' 9-461-" J2 IS NOT VALID HERE

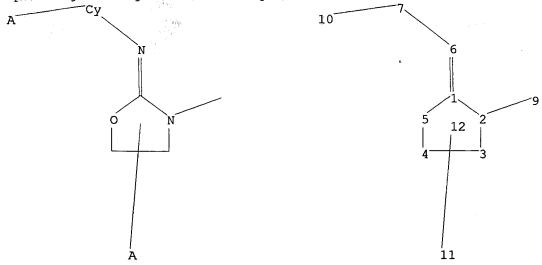
To display more answers, enter the number of answers you would like to see. To end the display, enter "NoVE", "N", "0", or "END". HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1): "0 S2' 9-861-" J\* 15 NOT VALID HERE

To display more answers, enter the number of answers you would like to see. To end the display, enter "NoNE", "N", "0", or "END". HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1): '0 SZ'  $\theta-\theta d1$ -" J\*' IS NOT VALID HERE

To display more answers, enter the number of answers you would like to see. To end the display, enter "MONE", "N", "0", or "END".
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? {1}:0

=>

Uploading C:\Program Files\Stnexp\Queries\10004396.str



chain nodes : 6 7 11 ring nodes : 1 2 3 4 5 ring/chain nodes : 9 10 chain bonds : 1-6 2-9 6-7 7-10 ring bonds : 1-2 1-5 2-3 3-4 4-5 exact/norm bonds : 1-2 1-6 2-3 2-9 6-7 exact bonds : 1-5 3-4 4-5 isolated ring systems : containing 1:

Match level:
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:Atom 9:CLASS 10:CLASS 11:CLASS 12:CLASS

#### L3 STRUCTURE UPLOADED

=> s 13

SAMPLE SEARCH INITIATED 19:08:22 FILE 'REGISTRY'
SAMPLE SCREEN SEARCH COMPLETED - 101 TO ITERATE

100.0% PROCESSED 101 ITERATIONS 10 ANSWERS SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*
PROJECTED ITERATIONS: 1418 TO 2622

PROJECTED ANSWERS: 11 TO

L4 10 SEA SSS SAM L3

=> d scan

L4 10 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Oxazolidine, 3-allyl-2-[(4-bromo-2-chlorophenyl)imino]- (8CI)
MF C12 H12 Br C1 N2 0

CH2-CH=CH2

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):2

L4 10 ANSWERS REGISTRY COPYRIGHT 2003 ACS IN Oxazolidine, 3-allyl-2-[(2,4-dichlorophenyl)imino]- (8CI) MF C12 H12 C12 N2 O

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

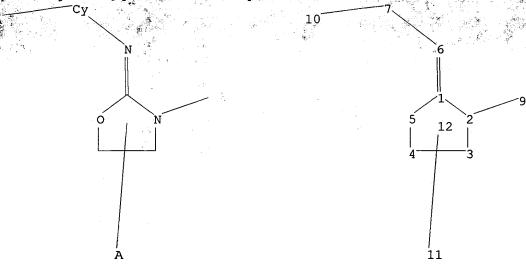
L4 10 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzenamine, N-[4-ethenyl-3-(2-methylphenyl)-2-oxazolidinylidene]-2-methyl(5)- (9CI)
MF C19 H20 N2 O

Absolute stereochemistry.
Double bond geometry unknown.

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

Uploading C:\Program Files\Stnexp\Queries\10004396.str



chain nodes : 6 7 11 ring nodes : 1 2 3 4 5 ring/chain nodes : 9 10 chain bonds : 1-6 2-9 6-7 7-10 ring bonds : 1-2 1-5 2-3 3-4 4-5 exact/norm bonds : 1-2 1-6 2-3 2-9 6-7 exact bonds : 1-5 3-4 4-5 isolated ring systems : containing 1:

Match level:
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:CLASS 7:Atom 9:CLASS 10:CLASS 11:CLASS 12:CLASS

# L5 STRUCTURE UPLOADED

=> s 15 SAMPLE SEARCH INITIATED 19:09:54 FILE 'REGISTRY' SAMPLE SCREEN SEARCH COMPLETED - 101 TO ITERATE

100.0% PROCESSED 101 ITERATIONS 2 ANSWERS SEARCH TIME: 00.00.01

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*

BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 1418 TO 2622

PROJECTED ANSWERS: 2 TO 124

L6 2 SEA SSS SAM L5

=> d scan

L6 2 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzenamine, N-[4-ethenyl-3-(2-methylphenyl)-2-oxazolidinylidene]-2-methyl, (s)- (9CI)
MF C19 H20 N2 O

Absolute stereochemistry. Double bond geometry unknown.

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):1

L6 2 ANSWERS REGISTRY COPYRIGHT 2003 ACS
IN Benzenamine, N-[5-ethyl-3-[3-(trifluoromethyl)phenyl]-2-oxazolidinylidene]3-fluoro- (9CI)
MF C18 H16 F4 N2 O
CI COM

\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

ALL ANSWERS HAVE BEEN SCANNED

=> s 15 sss full FULL SEARCH INITIATED 19:10:21 FILE 'REGISTRY' FULL SCREEN SEARCH COMPLETED - 1976 TO ITERATE

100.0% PROCESSED 1976 ITERATIONS

115 ANSWERS

SEARCH TIME: 00.00.01

L7 115 SEA SSS FUL L5

=> file ca caold COST IN U.S. DOLLARS

SINCE FILE TOTAL ENTRY SESSION

152.76 152.55 FULL ESTIMATED COST

FILE 'CA' ENTERED AT 19:10:55 ON 15 JAN 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 AMERICAN CHEMICAL SOCIETY (ACS)

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=> s 17

 $r_8$ 21 L7

=> sort py 18 SORT ENTIRE ANSWER SET? (Y)/N:.

2 ANSWERS DID NOT HAVE 'PY' SORT FIELD

PROCESSING COMPLETED FOR L8 L9 21 SORT L8 PY

=> d 19-21 cbib pi hitstr

L9 ANSWER 19 OF 21 CA COPYRIGHT 2003 ACS
136:216736 Preparation and was of: e.g., 2-arylimino-1,3-thiazolidines as progesterone receptor binding ligands, Dixon, Brian R., Bagi, Cedo M., Brennan, Catherine R., Britchli, David R., Bullock, William H., Chen, Jinshan, Collibes, William L.; Dally, Robert, Johnson, Jeffrey S., Kluender, Harold C. E.; Lathrop, William F., Liu, Pelyang Mass, Carol Ann. Redman, Aniko M., Scott, William J., Uchahns, Klaus; Wolanin, Donald Of (Bayer Corp., USA). U.S. US 6353006-Bl 20020305, 148 pp. (English). CODDN: USXANA. Replication. US 1999-453613 1991203. PRIORITY: US 6353006 Bl 20020305 US 1999-453613 1991203
II 285123-97-17 285123-88-89 285123-89-89
285123-90-27, 2-(2-Methyl-4-nitrophenylimino)-3-cyclopentyl-4, 4-dimethyl-1, 3-oxazolidine 285123-91-37, 2-(4-Cyano-2-ethylpenylimino)-3-cyclopentyl-4, 4-dimethyl-1, 3-oxazolidine RN: PAC (Pharmacological activity), SPN (Synthetic preparation), THU (Therapeutic use), BIOL (Biological study), PREP (Preparation), USES (Uses) (Uses)
(drug; preparation and use of, e.g., 2-arylimino-1;3-thiazolidines as progesterone receptor binding ligands)
285123-87-7 CA
Benzenamine, 2-methyl-N-[(45)-4-(1-methylethyl)-3-(2-methylpropyl)-2-oxazolidinylidene]-4-nitro- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.



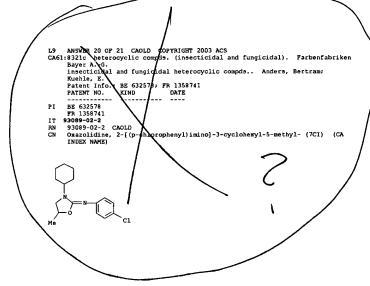
285123-88-8 CA Benzenamine, N-[(4S)-3,4-bis(2-methylpropyl)-2-oxazolidinylidene]-2-methyl-4-nitro-(SCI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

285123-89-9 CA
Benzonitrile, 4-[[(4S)-3,4-bis(2-methylpropyl)-2-oxazolidinylidene]amino]-3-ethyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

Double bond geometry unknown.



CA, COPYRIGHT 2003 ACS ANSWER 19 OF 21 (Continued)

285123-90-2 CA Benzehamine, N-i3-cyclopentyl-4,4-dimethyl-2-oxazolidinylidene)-2-methyl-4 nitro- (SCI) (CA INDEX NAME)

285123-91-3 CA
Benzonitrile, 4-[(3-cyclopentyl-4,4-dimethyl-2-oxazolidinylidene)amino]-3-ethyl- (9C1) (CA INDEX NAME)

L9 ANSWER 21 OF 21 CAOLD COPYRIGHT 2003 ACS
CA61:5637h acetylenic amines - (VIII) cyclization of acetylenic ureas.
Easton, Nelson R.; Cassady, D. R.; Dillard, R. D.
IT 93187-17-8 CAOLD
RN 93187-17-8 CAOLD
CN OXACOLdidine, 2-[(p-chlorophenyl)imino]-3,4,4-trimethyl-5-methylene- (7CI) (CA INDEX NAME)

=> file reg
COST IN U.S. DOLLARS

FULL ESTIMATED COST

SINCE FILE TOTAL ENTRY SESSION 9.82 162.58

FILE 'REGISTRY' ENTERED AT 19:13:07 ON 15 JAN 2003 USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT. PLEASE SEE "HELP USAGETERMS" FOR DETAILS. COPYRIGHT (C) 2003 American Chemical Society (ACS)

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STRUCTURE FILE UPDATES: 14 JAN 2003 HIGHEST RN 479024-64-1 DICTIONARY FILE UPDATES: 14 JAN 2003 HIGHEST RN 479024-64-1

TSCA INFORMATION NOW CURRENT THROUGH MAY 20, 2002

Please note that search-term pricing does apply when conducting SmartSELECT searches.

Crossover limits have been increased. See HELP CROSSOVER for details.

Experimental and calculated property data are now available. See HELP PROPERTIES for more information. See STNote 27, Searching Properties in the CAS Registry File, for complete details: http://www.cas.org/ONLINE/STN/STNOTES/stnotes27.pdf

=> s pyridine L10 582266 PYRIDINE

=> d rid

'RID' IS NOT A VALID FORMAT FOR FILE 'REGISTRY'

The following are valid formats:

Substance information can be displayed by requesting individual fields or predefined formats. The predefined substance formats are: (RN = CAS Registry Number)

REG - RN

SAM - Index Name, MF, and structure - no RN FIDE - All substance data, except sequence data

IDE - FIDE, but only 50 names
SQIDE - IDE, plus sequence data

SQIDE3 - Same as SQIDE, but 3-letter amino acid codes are used

SQD - Protein sequence data, includes RN

SQD3 - Same as SQD, but 3-letter amino acid codes are used

SQN - Protein sequence name information, includes RN

CALC - Table of calculated properties EPROP - Table of experimental properties

PROP - EPROP and CALC

Any CA File format may be combined with any substance format to obtain CA references citing the substance. The substance formats must be cited first. The CA File predefined formats are:

ABS -- Abstract

APPS - Application and Priority Information

BIB -- CA Accession Number, plus Bibliographic Data

CAN -- CA Accession Number

CBIB -- CA Accession Number, plus Bibliographic Data (compressed)

IND -- Index Data

IPC -- International Patent Classification

PATS -- PI, SO

STD -- BIB, IPC, and NCL

IABS --ABS, indented, with text labels

IBIB -- BIB, indented, with text labels

ISTD -- STD format, indented

OBIB ----- AN, plus Bibliographic Data (original)

OIBIB ----- OBIB, indented with text labels

SBIB ----- BIB, no citations

SIBIB ----- IBIB, no citations

The ALL format gives FIDE BIB ABS IND RE, plus sequence data when it is available.

The MAX format is the same as ALL.

The IALL format is the same as ALL with BIB ABS and IND indented, with text labels.

For additional information, please consult the following help messages:

HELP DFIELDS -- To see a complete list of individual display fields. HELP FORMATS -- To see detailed descriptions of the predefined formats. ENTER DISPLAY FORMAT (IDE):rsd

# L10 ANSWER 1 OF 582266 REGISTRY COPYRIGHT 2003 ACS

Ring System Data

Analysis EA	Sequence	e ithe R	ings  Form	RID	ier Occurrence
C6 C5N	IC6	16	1C6 1C5N	46.150.  46.156.	18   1

=> s pyridine/cn L11 1 PYRIDINE/CN => d rsd

| Lil | AMSWER | 1 OF 1 REGISTRY COPYRIGHT 2003 ACS | Ring System Data | Elemental| Elemental| Size of | Ring System | Ring | RID | Analysis | Sequence | the Rings| | Formula | | Identifier | Occurrence | EA | ES | SZ | RE | | RID | Count

L1

(FILE 'HOME' ENTERED AT 19:03:27 ON 15 JAN 2003)

FILE 'REGISTRY' ENTERED AT 19:03:34 ON 15 JAN 2003 STRUCTURE UPLOADED

L2 15 S L1

L3 STRUCTURE UPLOADED

L4 10 S L3

L5 STRUCTURE UPLOADED

L6 2 S L5

L7 115 S L5 SSS FULL

FILE 'CA, CAOLD' ENTERED AT 19:10:55 ON 15 JAN 2003

L8 21 S L7

L9 21 SORT L8 PY

FILE 'REGISTRY' ENTERED AT 19:13:07 ON 15 JAN 2003

L10 582266 S PYRIDINE

L11 1 S PYRIDINE/CN

=> s 17 and 46.156/rid

1552842 46.156/RID

L12 0 L7 AND 46.156/RID

=> s 17 and C5N/rf

1557529 C5N/RF

L13 0 L7 AND C5N/RF

=> file ca

COST IN U.S. DOLLARS SINCE FILE

ENTRY SESSION 19.60 182.18

TOTAL

FULL ESTIMATED COST

FILE 'CA' ENTERED AT 19:16:18 ON 15 JAN 2003
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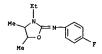
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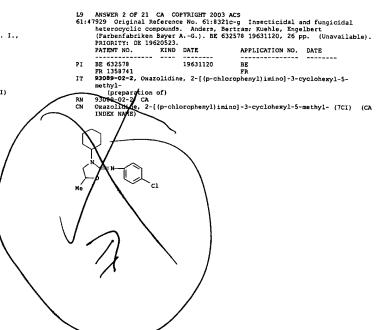
FILE COVERS 1907 - 9 Jan 2003 VOL 138 ISS 3 FILE LAST UPDATED: 9 Jan 2003 (20030109/ED)

This file contains CAS Registry Numbers for easy and accurate substance identification.

=> d 19 cbib pi fhitstr 1-10
YOU HAVE REQUESTED DATA FROM FILE 'CA, CAOLD' - CONTINUE? (Y)/N:.

US 2902356 1959 2260-79-9 CAOLD Oxazolidine, 3-ethyl-2-{(p-fluorophenyl)imino]-4,5-dimethyl- (6CI, 8CI) (CA INDEX NAME)





L9 ANSWER 3 OF 21 CA COPYRIGHT 2003 ACS
61:32397 Original Reference No. 61:5637h,5638a-b Acetylenic amines. VIII.
Cyclization of acetylenic ureas. Easton, Nelson R.; Cassady, Donald R.;
Dillard, Robert D. (Eii Lilly and Co., Indianapolis, IN). J. Org. Chem.,
29(7), 1851-5 (Unavailable) 1964. CODEN: JOCEAH. ISSN: 0022-3263.

IT 93187-17-8, Oxazolidine, 2-[(p-chlorophenyl)imino]-3,4,4-trimethyl5-methylene(preparation of)
RN 93187-17-8 CA
Oxazolidine, 2-((p-chlorophenyl)imino]-3,4,4-trimethyl-5-methylene- (7CI)
(CA INDEX NAME)

L9 ANSWER 4 OF 21 CA COPYRIGHT 2003 ACS
71:124419 Thiazole derivatives with bactericidal activity. Kutter, Eberhard;
Reuter, Wolfgang, Sauter, Roberty Machleidt, Hansy Wildfewer, Alexander;
Goeth, Hanns (Thomas, Dr. Karl, G.m.b. H.). S. African ZA 6807440
19690423, 36 pp. (Russian). CODEN: SFXXAB. PRIORITY: DE 19671116 19681001 19681001.
PATENT NO. KIND DATE APPLICATION NO. DATE

PI ZA 6807440 19690423 ZA 6807440 19690423
24240-64-OP
RL: SPN (Synthetic preparation); PREP (Preparation) (preparation of)
24240-64-O CA
3-Oxazoligineethanol, a,5-dimethyl-2-[(5-nitro-2-thiazolyl)imino]-(8CI) (GA INDEX NAME) Mé NO2

L9 ANSWER 5 OF 21 CA COPYRIGHT 2003 ACS
74:31746 Antibacterial 2-amino-5-nitrothiazoles substituted on the amino group. (Thomae, Dr. Karl, G.m.b.H.). fr. FR 1592266 19700619, 17 pp. (French). CODEN: FRXXAX APPLICATION FR 1968-1592266 19681115.
PATENT NO. KIND DATE APPLICATION NO. DATE APPLICATION NO. DATE PR 1592266 A 19700719 PR 1968-159026 19681101
DE 1800074 A 19700709 DE 1968-1800074 19681001
BR 6804058 AO 19730703 DR 1968-204058 19681114 FR 1592266 A 19700511 FR 1968-15922 DE 1800074 A 19700709 DE 1968-18000 BR 6804058 AO 19730703 BR 1968-20405 24240-64-0 CA CPE (preparation of) 24240-64-0 CA 3-Oxazolidineethanol, a5-dimeth/1-2-[(5-nitro-2-th/68CI) (CA INDEX NAME) IT 1-2-[(5-nitro-2-thiazolyl)imino]-

L9 ANSWER 6 OF 21 CA COPYRIGHT 2003 ACS

105:114951 Synthesis of 2-(2-arylimino-3-alkyl-4-oxazolidinyl)-Nalkylacetamides and 2-(2-arylimino-3-alkyl-4-thiazolidinyl)-Nalkylacetamides from 2(5H)-furanone. Tyukhteneva, Z. I., Badovskaya, L.
A.; Kozlovskaya, I. N.; Muzychenko, G. F. (Politekh. Inst., Krasnodar,
350006, USSR). Khimiya Geterotiskilcheskikh Seedinenii (12), 1629-32
(Russian) 1985. CODEN: KGSSAQ. ISSN: 0453-8274. OTHER SOURCES: CASREACT
105:114951.

IT 104053-32-9P
RL: SPN (Synthetic preparation); PREP (Preparation)
(preparation of)
RN 104053-32-9 CA
4-0xazolidineacetamide, N,3-dibutyl-2-[(4-chlorophenyl)imino]- (9CI) (CA
INDEX NAME) 0 || n-BuNH-C-CH2 `c1

L9 ANSWER 7 OF 21 CA COPYRIGHT 2003 ACS
111:7388 Process for the preparation of iminooxazolidines useful as herbicides. Felix, Raymond Anthony (Stauffer Chemical Co., USA). Braz. Pedido PI BR 8705543 A 19880524, 48 pp. (Portuguese). CODEN: BPXXDX. APPLICATION: BR 1987-5543 19871016. PRIORITY: US 1986-92014 19861017; US 1986-92014 3 19861017.

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	BR 8705543	Α	19880524	BR 1987-5543	19871016
	US 4806653	Α	19890221	US 1986-920345	19861017
	ES 2036583	T3	19930601	ES 1987-309069	19871014
	JP 63152369	A2	19880624	JP 1987-258481	19871015
	CA 1328455	A1	19940412	CA 1987-549329	19871015
	DK 9705408	A	19880418	DK 1987-5408	19871016
	AU 8779852	A1	19880428	AU 1987-79852	19871016
	AU 592233	B2	19900104		
	HU 44906	A2	19880530	HU 1987-4667	19871016
	CN '87107006	A	19880629	CN 1987-107006	19871016
	CN 1017996	В	19920826		
	ZA 8707785	A	19880831	ZA 1987-7785	19871016
	IN 165720	A	19891223	IN 1987-MA745	19871016
	SU 1776185	A3	19921115	SU 1987-4203487	19871016
	SU 1681726	A3	19910930	SU 1988-4355057	19880201
	US 4900351	A	19900213	US 1988-167934	19880314

US 4900351 A 19900213 US 1988-107534 1555551
IT 115315-46-3p
RL: AGR (Agricultural use); BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); BIOL (Biological study); PREP (Preparation); USES (Uses) (preparation of, as herbicide)
RN 115315-46-3 CA CN
Benzenamine, 4-chloro-N-[5-ethyl-3-[3-(trifluoromethyl)phenyl]-2-oxazolidinylidene]- (9CI) (CA INDEX NAME)

L9 ANSWER 8 OF 21 CA COPYRIGHT 2003 ACS
109:230998 Preparation of owazolidines and thiazolidines as acaricides.
15hintsu, Keitchi; Imagawa, Hiroyuki; Yamada, Tomio; Matsuda, Michihiko;
Kitagawa, Yukio (Nippon Soda Co., Ltd., Japan). Upn. Kokai Tokkyo Koho JP
63041471 Az 1980222 Showa, 10 pp. (Japanese). CODEN: JXXXAF.
APPLICATION: JP 1986-186454 19860808.
APPLICATION NO. DATE

NO. TREASURE AND DATE APPLICATION NO. DATE JP 63041471 A2 19880222 JP 1986-186454 19860808
117518-52-2P
RL: SPN (Synthetic preparation), PREP (Preparation)
(preparation on as acaricide)
117518-52-2 CA
3-Oxazolidinec/srboxamide, 2-[(4-chloro-2-methylphenyl)imino]-N-cyclohexyl-4-methyl-5-(4/methylphenyl)- (9CI) (CA INDEX NAME)

L9 ANSWER 9 OF 21 CA COPYRIGHT 2003 ACS
109:54765 Preparation and testing of diaryliminooxazolidines as pre- and postemergent herbicides. Felix, Raymond Anthony (Stauffer Chemical Co., USA). Eur. Pat. Appl. EP 265162 Al 1980427, 22 pp. DESIGNATED STATES: R: AT. BE, CH, DE, ES, FR, GB, GR, IT, LI, NL, SE. (English). CODEN: EPXKDW. APPLICATION: EP 1897-309069 19871014. PRIORITY: US 1986-920014 19861017. US 1986-920345 19861017. APPLICATION NO. DATE

	PA'	TENT NO.		KIND	DATE	APPLICATION NO.	DATE
21	EP	265162		A1		EP 1987-309069	19871014
	EP	265162		B1	19910306		
		R: AT,	BE,	CH, DE	, ES, FR, G	B, GR, IT, LI, NL, SE	3
	US	4806653		A	19890221	US 1986-920345	19861017
	AT	61358		E	19910315	AT 1987-309069	19871014
	ES	2036583		т3	19930601	ES 1987-309069	19871014
	JP	63152369		A2	19880624	JP 1987-258481	19871015
	CA	1328455		A1	19940412	CA 1987-549329	19871015
	DK	8705408		A	19880418	DK 1987-5408	19871016
	AU	8779852		A1	19880428	AU 1987-79852	19871016
	ΑU	592233		B2	19900104		
	HU	44906		A2	19880530	HU 1987-4667	19871016
	CN	87107006		A	19880629	CN 1987-107006	19871016
	CN	1017996		В	19920826		
	ZA	8707785		A	19880831	ZA 1987-7785	19871016
	IN	165720		A	19891223	IN 1987-MA745	19871016
	SU	1776185		A3	19921115	SU 1987-4203487	19871016
	SU	1681726		A3	19910930	SU 1988-4355057	19880201
	US	4900351		A	19900213	US 1988-167934	19880314

US 490031 A 19900213 US 1988-167934 19880314
118318-46-3P A 19900213 US 1988-167934 19880314
118315-46-3 CA 19900213 US 1988-167934 19880314
118315-46-3P A 19900214
118315-46-3P A 19880314
118315-46-3P A 19900214
118315-46

19 ANSWER 10 OF 21 CA COPYRIGHT 2003 ACS
126:293283 Palladium(0)-Catalyzed Asymmetric Cycloaddition of Vinyloxiranes
with Heterocumulenes Using Chiral Phosphine Ligands: An Effective Route to
Highly Enantioselective Vinyloxazolidine Derivatives. Larksarp,
Chitchamair Alper, Howard (Department of Chemistry, University of Ottawa,
Ottawa, ON, KIN 6W5, Can.). Journal of the American Chemical Society,
119 [16], 3709-3715 (English) 1997. CODEN: JACSAT. ISSN: 0002-7863.

IT 189128-42-5 FR.
RI: SPM (Synthetic preparation), PREP (Preparation)
(palladium(0)-catalyzed asym. cycloaddn. of vinyloxiranes with
heterocumulenes using chiral phosphine ligands)
RN 189128-42-5 CA
CN Benzenamine, 4-chloro-N-[3-(4-chlorophenyl)-4-ethenyl-2-oxazolidinylidene](9CI) (CA INDEX NAME)

$$H_2C = CH$$
 $N$ 
 $N$ 
 $C1$ 

=> d 19 cbib pi fhitstr 11-18
YOU HAVE REQUESTED DATA FROM FILE 'CA, CAOLD' - CONTINUE? (Y)/N:.

129:230662 Highly Enanticoslective Synthesis of 1,3-Oxazolidin-2-imine
Derivatives by Asymmetric Cycloaddition Reactions of Vinyloxiranes with
Unsymmetrical Carbodiimides Catalyzed by Palladium(0) Complexes,
Larksarp, Chitchamair Alper, Howard (Department of Chemistry, University
of Ottawa, Ottawa, ON, KIN SNS, Can.), Journal of Organic Chemistry,
63(18), 6229-6233 (English) 1998. CODEN: JOCEAH. ISSN: 0022-3263. OTHER
SOURCES: CASREACT 129:230662. Publisher: American Chemical Society.
RI: SPN (Synthetic preparation); PREP (Preparation)

212697-18-59
RE: SPN (Synthetic preparation); PREP (Preparation)
(stereoselective preparation of oxazolidinimines by asym. cycloaddn. of vinyloxicanes with unsym. carbodiimides)
212697-15-5 CA
Benzenamine, N-(3-cyclohexyl-4-ethenyl-2-oxazolidinylidene)-4-methoxy-(9CI) (CA INDEX NAME)

L9 ANSWER 13 OF 21 CA COPYRIGHT 2003 ACS

133:120323 Preparation of 2-aryliminothiazolidines and related compds. progesterone receptor binding agents. Dixon, Brian R.; Bagi, Cedo M.; Brennan, Catherine R.; Beittelli, David R.; Bullock, William H.; Chen, Jinshan; Collibee, William L.; Dally, Robert Johnson, Jeffrey S.; Kluender, Harold C. E.; Lathrop, William F.; Liu, Peiying; Mase, Carol Ann; Redman, Aniko M.; Scott, William J.; Urbahns, Klaus; Wolanin, John (Bayer Corporation, USA). PCT Int. Appl. WO 2000042031 A2 20000720, 274 pp. DESIGNATED STATES: W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CM, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, KN, NO, NZ, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, RG: AT, BE, BF, BJ, CF, CG, CH, CT, CM, CY, DE, DK, ES, FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG. (English). CODDEY: PIXXDIS. APPLICATION: WO 1999-US29601 19991214.

PATEMIT NO. KIND DATE

PI WO 2000042031 A2 20000720 WO 1999-US29601 19991214

PATEMIT NO. KIND DATE

APPLICATION NO. DATE

PI WO 2000042031 A2 20000720 WO 1999-US29601 19991214

PATEMIT NO. KIND DATE

APPLICATION NO. DATE

PI WO 2000042031 A2 20000720 WO 1999-US29601 19991214

PATEMIT NO. KIND DATE

APPLICATION NO. DATE

PI WO 2000042031 A2 20000720 WO 1999-US29601 19991214

PATEMIT NO. KIND DATE

PI WO 2000042031 A2 20000720 WO 1999-US29601 19991214

PATEMIT NO. KIND DATE

PI WO 2000042031 A3 2001103

PI AR, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU, CZ, OE, DK, DM, KE, ES, FI, GB, GD, GE, GH, GH, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MK, MO, NZ, PL, PT, PO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BA, BB, BG, BR, IT, LU, W, N, Y, YU, ZA, ZW, AM, AZ, BA, BB, BG, R

285123-97-79
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified), SPN (Synthetic preparation); THU (Therapeutic use); BIOL (Biological study); PREF (Preparation); USES (Uses) (preparation of 2-aryliminothiazolidines and related compds. progesterone receptor binding agents)
285123-97-7 CA
Benzenamine, 2-methyl-N-[(4S)-4-(1-methylethyl)-3-(2-methylpropyl)-2-oxazolidinylidene]-4-nitro- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

L9 ANSWER 12 OF 21 CA COPYRIGHT 2003 ACS
128:217324 The synthesis of 2-imidazolidones on solid support by tandem
aminoacylation/Michael addition. Goff, Dane (Chiron Corp., Emeryville,
CA, 94608, USA). Tetrahedron Letters, 39(12), 1477-1480 (English) 1998.
CODEN: TELEAY. ISSN: 0040-4039. Publisher: Elsevier Science Ltd..
IT 204330-13-2P

ZVes30-13-2F
RL: SPN (Synthetic preparation): PREP (Preparation)
 (preparation of imidazolidones on solid support by tandem aminoacylation/Michael addition)
204330-13-2 CA

ביי בייטין ב- CA 5-Owazolidineacetamide, 2-[(2-ethoxyphenyl)imino]-3-(2-methylpropyl)-(9CI) (CA INDEX NAME)

132:222542 Preparation of 1-(3-heterocyclylphenyl)isothioureas, -isoureas, -guanidines and -amidines as herbicides. Karp, Gary Mitchell (American Cyanamid Company, USA). Bur. Pat. Appl. EP 985670 Al 20000315, 303 pp. DESIGNATED STATES: R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO. (English). CODEN: EFXKUW. APPLICATION: EP 1999-306382 19990812. PRICRITY: US 1998-133872 19980813. PATENT NO. KIND DATE APPLICATION NO. DATE

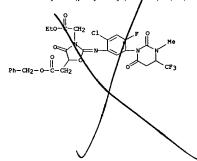
PI EP 985670 Al 20000315 EP 1999-306382 19990812
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO

17 260977-27-3P
RL: AGR (Agricultural use), BAC (Biological activity or effector accept

IT 260977-27-3P

RL: AGR (Agricultural use), BAC (Biological activity or effector, except adverse): BSU (Biological study, unclassified): SFN (Synthetic preparation): BIOL (Biological study): PREP (Preparation): USES (Uses) (preparation of 1-(3-heterocyclylphenyl)isothioureas, -isoureas, -guanidines

(preparation of 1-(3-heterocyclylphenyl)isothioureas, -isoureas,
nidines
and -amidines as herbicides)
260977-27-3 CA
3,5-Oxazolidinediacetic acid, 2/[[2-chloro-4-fluoro-5-[tetrahydro-3-methyl-2,6-dioxo-4-(trifluoromethyl)-/(ZE)-pyrimidinyl)phenyl]imino]-4-oxo-,
α3-ethyl α5-(phenylmethyl) esper (9CI) (CA INDEX NAME)



L9 ANSWER 15 OF 21 CA COPYRIGHT 2003 ACS
137:33286 Fluoric derivative of N-phenyl-2-oxazolylamino preparing process and
its application. Qian, Xuhong; Song, Gonghua; Li, Zhibin; Yu, Hongzhong;
Li, Zhong; Liu, Zhi (Huadeng Univ. of Science and Engineering, Peop. Rep.
China). Faming Zhuanli Shenqing Gongkai Shuomingshu CN 1294126 A
20010509, 16 pp. (Chinese). CODEN: CNXXEV. APPLICATION: CN 2000-127297
20001107.
PATENT NO. XIMD DATE
APPLICATION NO. DATE CN 2000-127297 20001107 20010509

CN 1294126
436854-67-OP
RL: AGR (Agricultus)
(Synthetic prepara
(Uses) ral use); BSU (Mological study, unclassified); SPN kion); BIOL Biological study); PREP (Preparation); USES

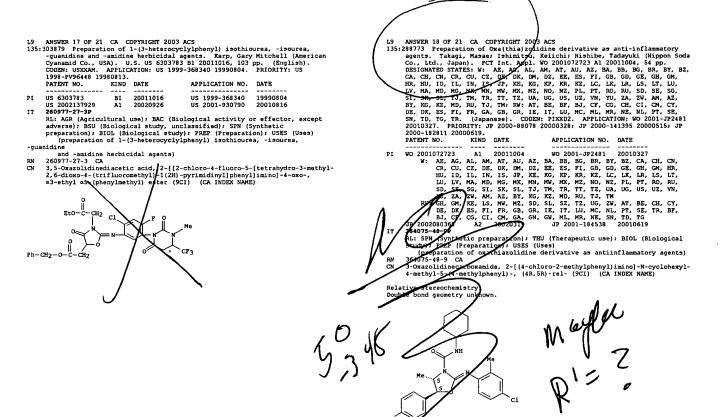
(Uses)
(prepn of fluoric derivative of N-phenyl-2-oxazolylamino and its application)
436854-67-0 CB.
1,2,3,4-Butanetetrol, 1-[2-[(4-hydroxyphenyl)imino]-3-methyl-5-oxazolidinyl]- (9CI) (CA INDEX NAME)

L9 ANSWER 16 OF 21 CA COPYRIGHT 2003 ACS 136:247793 Syntheses of 2-aryliminooxazolidine derivatives as trehalase inhibitors. Qian, Xuhong, Li, Zhibin, Liu, Zhir Song, Gonghuar Li, Zhong (State Key Laboratory of Fine Chemicals, Dalian University of Technology, Dalian, 116012, Peop. Rep. China). Journal of Antibiotics, 54(12), 1108-1110 (English) 2001. CODEN: JANTAJ. ISSN: 0021-8820. Publisher: Japan Antibiotics Research Association.
IT 404352-94-99

RI: BSU (Biological study, unclassified); SPN (Synthetic preparation);
BIOL (Biological study); PREP (Preparation)
(preparation of 2-aryliminooxazolidine derivs. as trehalase inhibitors

fungicides) 404352-94-9 CA 1,2,3,4-8utanetetrol, 1-[(55)-2-[(4-hydroxyphenyl)imino]-3-methyl-5-oxazolidinyl]-, (15,2R,3R)- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.



=> s 17 and 46.156/rid
'RID' IS NOT A VALID FIELD CODE

18 L7

0 46.156/RID

L14

0 L7 AND 46.156/RID

=> d 19 cbib pi hitstr 5 6 11-13
YOU HAVE REQUESTED DATA FROM FILE 'CA, CAOLD' - CONTINUE? (Y)/N:.

L9 ANSWER 5 OF 21 CA COPYRIGHT 2003 ACS
74:31746 Antibacterial 2-amino-5-nitrothiazoles substituted on the amino group. (Thomae, Dr. Karl, G.m.b.H.). Fr. FR 1592266 19700619, 17 pp. (French). CODEN: FRXXAK. APPLICATION: FR 1968-1592266 19681115.

PATENT NO. KIND DATE APPLICATION NO. DATE

FR 1592266 A 19700511 FR 1968-1592266 19681115
DE 1800074 A 19700709 DE 1968-1800074 19681001
BR 6804058 A0 19730703 DR 1968-204058 19681114 FR 1592266 A 19700511 FR 1968-1592266 19681115
DE 1800074 A 19700709 DE 1968-1800074 19681001
BR 6804058 AO 19730703 BR 1968-204058 19681014
24240-64-0P
RL: SPN (Synthetic preparation), PREP (Preparation)
(preparation of)
24240-64-0 CA
3-Oxazolidinethanol, α,5-dimethyl-2-[(5-nitro-2-thiazolyl)imino](8CI) (CA INDEX NAME)

L9 ANSWER 11 OF 21 CA COPYRIGHT 2003 ACS
129:230662 Highly Enantioselective Synthesis of 1,3-Oxazolidin-2-imine
Derivatives by Asymmetric Cycloaddition Reactions of Vinyloxiranes with
Unsymmetrical Carbodlimides Catalyzed by Palladium(0) Complexes.
Larksarp, Chitchamair Alper, Howard (Department of Chemistry, University
of Ottawa, Ottawa, ON, KIN GNS, Can.). Journal of Organic Chemistry,
63(18), 6229-623 (English) 1998. CODEM: JOCKEM. ISSN: 0022-3263. OTHER
SOURCES: CASREACT 129:230662. Publisher: American Chemical Society.
17 212687-13-59 212687-38-9 212687-26-5P
212687-22-49 212687-30-49 212687-36-5P
212687-38-99 212687-36-9 212687-37-1P
212687-38-99 212687-36-99 212687-37-1P
212687-38-99 212687-36-0P

212667-38-2P
RE: SPN (Synthetic preparation); PREP (Preparation)
(stereoselective preparation of oxazolidinimines by asym. cycloaddn. of
vinyloxiranes with unsym. carbodimides)
212687-15-5 CA
Benzenamine, N-(3-cyclohexyl-4-ethenyl-2-oxazolidinylidene)-4-methoxy(9CI) (CA INDEX NAME)

212697-18-8 CA
Benzenamine, N-(3-butyl-4-ethenyl-2-oxazolidinylidene)-4-chloro- (9CI)
(CA INDEX NAME)

$$\text{H}_2\text{C} = \text{CH} \bigvee_{N=-N}^{\text{n-Bu}} \text{C1}$$

212687-20-2 CA Benzenamine, N-(3-cyclohexyl-4-ethenyl-2-oxazolidinylidene)-4-fluoro-(9C1) (CA INDEX NAME)

212687-22-4 CA
Benzenamine, N-(3-cyclohexyl-4-ethenyl-2-oxazolidinylidene)-2,6-dimethyl-(9CI) (CA INDEX NAME)

ANSWER 11 OF 21 CA COPYRIGHT 2003 ACS

212687-24-6 CA Benzenamine, N-(4-ethenyl-3-ethyl-2-oxazolidinylidene)-4-methyl- [9CI] (CA INDEX NAME)

212687-26-8 CA
Benzenamine, N-(3-butyl-4-ethenyl-2-oxazolidinylidene)-2,6-dimethyl- (9CI)
(CA INDEX NAME)

212687-28-0 CA
Benzenamine, 4-chloro-N-(4-ethenyl-3-phenyl-2-oxazolidinylidene)- (9CI)
(CA INDEX NAME)

$$H_2C = CH$$

Ph

N

C1

212697-30-4 CA Benzenamine, N-(4-ethenyl-3-phenyl-2-oxazolidinylidene)-4-methyl- (9CI) (CA INDEX NAME)

$$H_2C = CH$$
 $N$ 
 $N$ 
 $Me$ 

212687-34-8 CA Benzenamine, N-[(45)-3-butyl-4-ethenyl-2-oxazolidinylidene]-4-chloro-

ANSWER 11 OF 21 CA COPYRIGHT 2003 ACS

212687-35-9 CA Forexament, N-[(45)-3-cyclohexyl-4-ethenyl-2-oxazolidinylidene]-4-fluoro-(951) [CA INDEX NAME]

Absolute stereochemistry. Double bond geometry unknown.

212687-36-0 CA
Benzenamine, N-[(4R)-3-cyclohexyl-4-ethenyl-2-oxazolidinylidene]-4-fluoro(9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

212687-37-1 CA
Benzenamine, N-[(4R)-3-cyclohexyl-4-ethenyl-2-oxazolidinylidene]-2,6dimethyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry. Double bond geometry unknown.

L9 ANSWER 12 OF 21 CA COPYRIGHT 2003 ACS

128:217324 The synthesis of 2-imidazolidones on solid support by tandem aminoacylation/Michael addition. Goff, Dane (Chiron Corp., Emeryville, CA, 94608, USA). Tetrahedron Letters, 39(12), 1477-1480 (English) 1998. CODEN: TELEAY. ISSN: 0040-4039. Publisher: Elsevier Science Ltd..

1204330-13-22 204330-14-3P

RL: SPN (Synthetic preparation): PREP (Preparation) (preparation of imidazolidones on solid support by tandem aminoacylation/Michael addition)

RN 204330-13-2 CA

CN 5-0Vascolidineacetamide, 2-[(2-ethoxyphenyl)imino]-3-(2-methylpropyl)-(9CI) (CA INDEX NAME)

204330-14-3 CA 5-0xazolidineacetamide, 2-[(2,4-dimethoxyphenyl)imino]-3-(2-methylpropyl)-(9CI) (CA INDEX NAME)

ANSWER 11 OF 21 CA 100

212687-38-2 CA Benzenamine, N-[(4R)-3-butyl-4-ethenyl-2-oxazolidinylidene]-2,6-dimethyl-(5C1) (CA INDEX NAME)

- 3

Absolute stereochemistry.
Double bond geometry unknown.

1.9 ANSWER 13 OF 21 CA COPYRIGHT 2003 ACS

133:120323 Preparation of 2-aryliminothiazolidines and related compds.

progesterone receptor binding agents. Dixon, Brian R.; Bagi, Cedo M.;
Brennan, Catherine R.; Brittelli, David R.; Bullock, William H.; Chen,
Jinshan; Collibee, William L.; Dally, Robert; Johnson, Jeffrey S.;
Kluender, Harold C. E.; Lathrop, William F.; Liu, Peiying; Mase, Carol
Ann; Redman, Aniko M.; Scott, William J.; Urbahns, Klaus; Wolanin, John J
(Bayer Corporation, USA). PCT Int. Appl. Wo 2000042031 A2 20000720, 274

pp. DESIGNATED STATES: W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY,
CA, CH, CN, CR, CU, CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR,
HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LK, LS, LT, LU, LV,
MA, MD, MG, MK, MM, MY, MX, NO, NZ, PL, PT, NO, RU, SD, SE, GG, SI, SK,
SL, TJ, TM, TR, TT, TZ, UA, UG, UZ, VN, YU, ZA, ZW, AM, AZ, BY, KG, KZ,
MD, RU, TJ, TH, RG: AT, BE, BF, BJ, CF, CG, CH, CH, CM, CV, DE, DK, ES,
FI, FR, GA, GB, GR, IE, IT, LU, MC, ML, MR, NE, NL, PT, SE, SN, TD, TG,
(English). CODEN: PIXXOZ APPLICATION: WO 1999-US29601 19991214.

PATEMT NO. KINO DATE

PI WO 2000042031 A2 20000720

WO 1999-US29601 19991214

WO 2000042031 A3 20001109

W: AE, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, CA, CH, CN, CR, CU,
CZ, DE, DK, DM, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL,
IN, 15, JF, RE, KG, KP, KR, KZ, LC, LK, LK, LS, LT, LU, LV, MA,
MD, MG, MK, MN, MW, MK, NO, XP, FL, FT, NO, RU, SD, SE, SG, SI, ST,
BY, KS, KZ, HD, RU, TJ, TM

RW: GH, GM, KM, LM, WK, NO, XP, FL, FT, NO, RU, SD, SE, SG, SI, ST,
MY, KS, KZ, HD, RU, TJ, TM

RW: GH, GM, KE, LS, MW, SD, SL, SZ, TZ, UG, ZW, AT, BE, CH, CY, DE,
DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, FT, SE, BF, BU, CF,
CG, CI, CM, GA, GN, GW, ML, MR, NE, SN, TD, TG

EP 1144386 A2 20011017

EP 19916999

A 200103318 A 2001003

NO 20010-3318 20010704

17 281523-90-72 285123-999-139

RL: BAC (Biological activity or effector, except adverse); BSU (Biological study); PREP (Preparation), 1885 (Hac-\*\*)

285123-90-2P 285123-91-3P
RL: BAC (Biological activity or effector, except adverse); BSU (Biological study, unclassified); SPN (Synthetic preparation); THU (Therapeutic use);
BIOL (Biological study); PREP (Preparation); USES (Uses)
(preparation of 2-aryliminothiazolidines and related compds. progesterone receptor binding agents)
285123-87-7 CA
Benzenamine, 2-methyl-N-[(4S)-4-(1-methylethyl)-3-(2-methylpropyl)-2-oxazolidinylidene]-4-nitro- (9CI) (CA INDEX NAME)

Absolute stereochemistry.
Double bond geometry unknown.

285123-88-8 CA
Benzenamine, N-[(45)-3,4-bis(2-methylpropyl)-2-oxazolidinylidene]-2-methyl-

Absolute stereochemistry.
Double bond geometry unknown.

RN 285123-89-9 CA
CN Benzonitrile, 4-[[(45)-3,4-bis(2-methylpropyl)-2-oxazolidinylidene]amino]3-ethyl- (9CI) (CA INDEX NAME)

Absolute stereochemistry.

RN 285123-90-2 CA
CN Benzenamine, N-(3-cyclopentyl-4,4-dimethyl-2-oxazolidinylidene)-2-methyl-4nitro- (9CI) (CA INDEX NAME)

RN 285123-91-3 CA
CN Benzonitrile, 4-[(3-cyclopentyl-4,4-dimethyl-2-oxazolidinylidene)amino]-3-ethyl- (9CI) (CA INDEX NAME)

=> logoff
ALL L# QUERIES AND ANSWER SETS ARE DELETED AT LOGOFF
LOGOFF? (Y)/N/HOLD:.

COST IN U.S. DOLLARS SINCE FILE TOTAL ENTRY SESSION

FULL ESTIMATED COST 2.00 259.96

STN INTERNATIONAL LOGOFF AT 19:27:05 ON 15 JAN 2003